

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page for STN Seminar Schedule - N. America
NEWS 2 AUG 15 CAOLD to be discontinued on December 31, 2008
NEWS 3 OCT 07 EPFULL enhanced with full implementation of EPC2000
NEWS 4 OCT 07 Multiple databases enhanced for more flexible patent
number searching
NEWS 5 OCT 22 Current-awareness alert (SDI) setup and editing
enhanced
NEWS 6 OCT 22 WPIDS, WPINDEX, and WPIX enhanced with Canadian PCT
Applications
NEWS 7 OCT 24 CHEMLIST enhanced with intermediate list of
pre-registered REACH substances
NEWS 8 NOV 21 CAS patent coverage to include exemplified prophetic
substances identified in English-, French-, German-,
and Japanese-language basic patents from 2004-present
NEWS 9 NOV 26 MARPAT enhanced with FSORT command
NEWS 10 NOV 26 MEDLINE year-end processing temporarily halts
availability of new fully-indexed citations
NEWS 11 NOV 26 CHEMSAFE now available on STN Easy
NEWS 12 NOV 26 Two new SET commands increase convenience of STN
searching
NEWS 13 DEC 01 ChemPort single article sales feature unavailable
NEWS 14 DEC 12 GBFULL now offers single source for full-text
coverage of complete UK patent families
NEWS 15 DEC 17 Fifty-one pharmaceutical ingredients added to PS

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that
specific topic.

All use of STN is subject to the provisions of the STN Customer
agreement. Please note that this agreement limits use to scientific
research. Use for software development or design or implementation
of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 10:14:38 ON 30 DEC 2008

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.42	0.42

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 10:15:31 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7
DICTIONARY FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

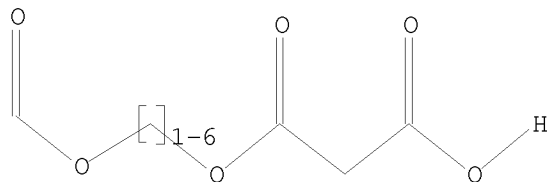
Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary
files\10531382\10531382 carbonyloxyalkyl esters.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam

SAMPLE SEARCH INITIATED 10:16:33 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 1768 TO ITERATE

100.0% PROCESSED 1768 ITERATIONS

2 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

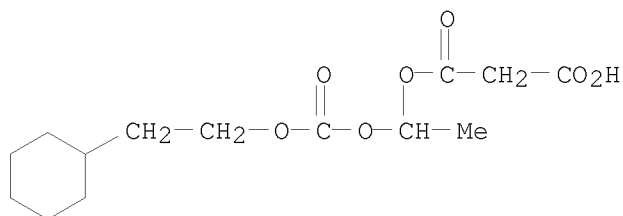
PROJECTED ITERATIONS: 32838 TO 37882

PROJECTED ANSWERS: 2 TO 124

L2 2 SEA SSS SAM L1

=> d scan

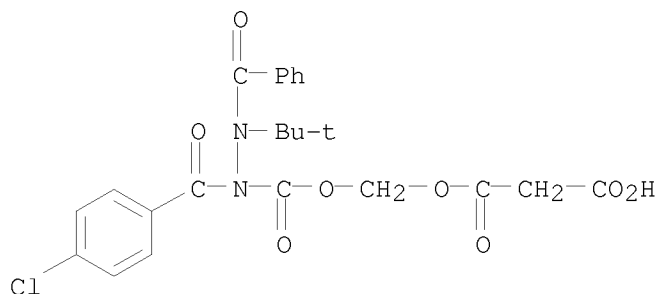
L2 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[1-[[2-cyclohexylethoxy)carbonyl]oxy]ethyl] ester
MF C14 H22 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L2 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[[[2-benzoyl-1-(4-chlorobenzoyl)-2-(1,1-dimethylethyl)hydrazinyl]carbonyl]oxy]methyl] ester
MF C23 H23 Cl N2 O8
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

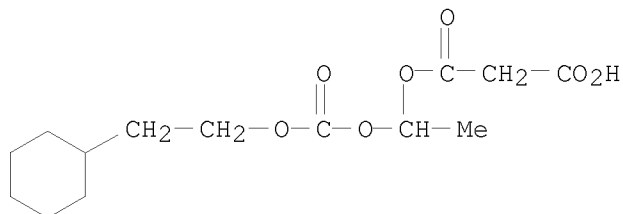
=> search l1 sss full
FULL SEARCH INITIATED 10:17:23 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 34586 TO ITERATE

100.0% PROCESSED 34586 ITERATIONS 40 ANSWERS
SEARCH TIME: 00.00.01

L3 40 SEA SSS FUL L1

=> d scan

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[1-[[2-cyclohexylethoxy)carbonyl]oxy]ethyl] ester
MF C14 H22 O7

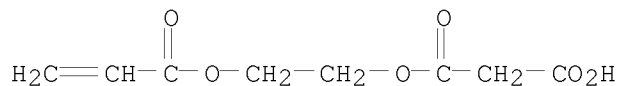


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):40

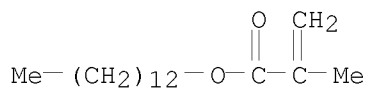
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 2-Propenoic acid, 2-methyl-, tridecyl ester, polymer with dodecyl
2-propenoate and 2-hydroxyethyl 2-propenoate,
2-[(1-oxo-2-propenyl)oxy]ethyl propanedioate (9CI)
MF (C17 H32 O2 . C15 H28 O2 . C5 H8 O3)x . x C8 H10 O6

CM 1

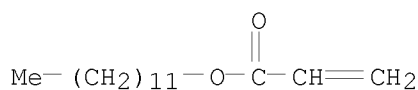


CM 2

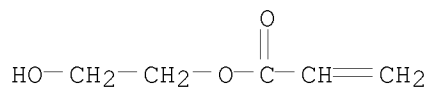
CM 3



CM 4

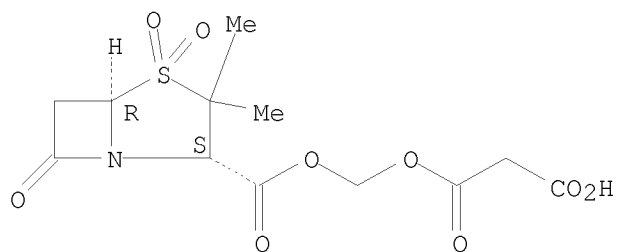


CM 5



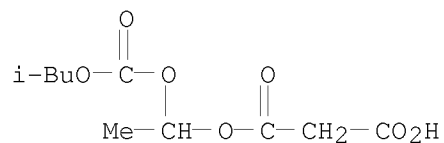
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, sodium salt, (2S-cis)- (9CI)
MF C12 H15 N O9 S . Na

Absolute stereochemistry.



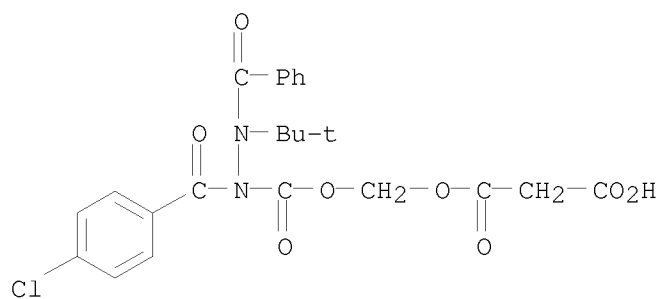
● Na

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[1-[(2-methylpropoxy)carbonyl]oxy]ethyl] ester
MF C10 H16 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

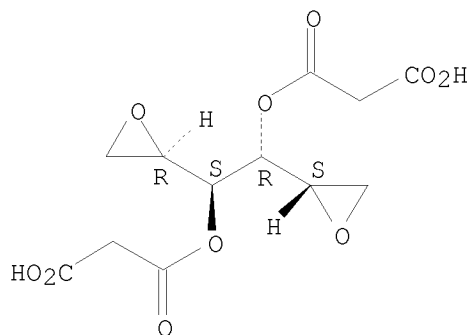
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[[[2-benzoyl-1-(4-chlorobenzoyl)-2-(1,1-dimethylethyl)hydrazinyl]carbonyl]oxy]methyl] ester
MF C23 H23 Cl N2 O8
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

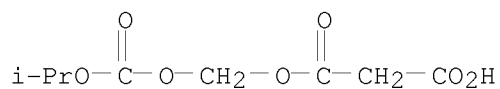
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Galactitol, 1,2:5,6-dianhydro-, bis(hydrogen propanedioate) (9CI)
 MF C12 H14 O10

Relative stereochemistry.



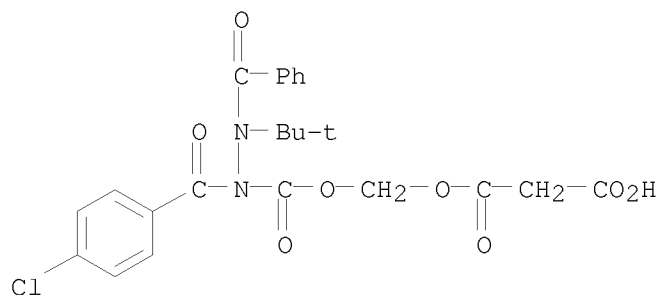
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[(1-methylethoxy)carbonyl]oxy]methyl] ester
 MF C8 H12 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

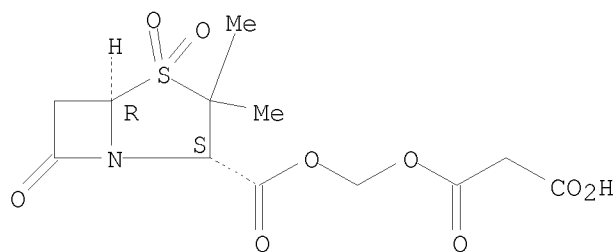
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[2-benzoyl-1-(4-chlorobenzoyl)-2-(1,1-dimethylethyl)hydrazinyl]carbonyl]oxy]methyl] ester, sodium salt (1:1)
 MF C23 H23 Cl N2 O8 . Na



● Na

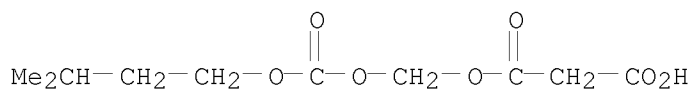
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, (2S-cis)- (9CI)
 MF C12 H15 N O9 S
 CI COM

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

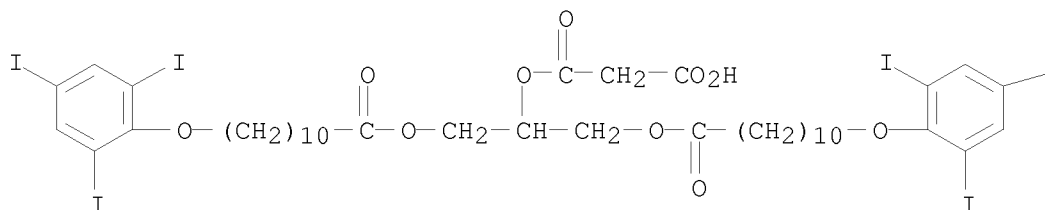
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[(3-methylbutoxy)carbonyl]oxy]methyl] ester
 MF C10 H16 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[2-[[1-oxo-11-(2,4,6-triiodophenoxy)undecyl]oxy]-1-
MF [[1-oxo-11-(2,4,6-triiodophenoxy)undecyl]oxy)methyl]ethyl] ester
C40 H52 I6 O10

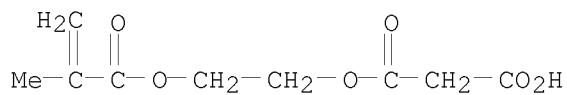
PAGE 1-A



PAGE 1-B

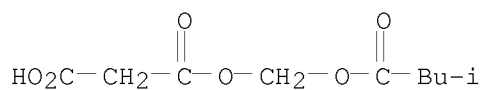
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[2-[(2-methyl-1-oxo-2-propen-1-yl)oxy]ethyl] ester
MF C9 H12 O6
CI COM



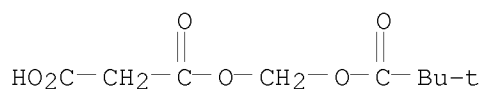
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[(3-methyl-1-oxobutoxy)methyl] ester
MF C9 H14 O6



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

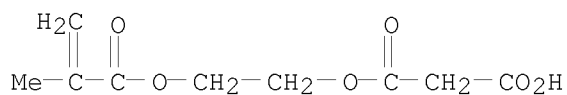
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[(2,2-dimethyl-1-oxopropoxy)methyl] ester
MF C9 H14 O6



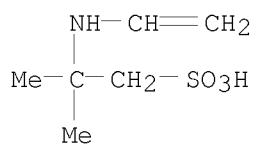
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, mono[2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl] ester,
polymer with 2-(ethenylamino)-2-methyl-1-propanesulfonic acid and methyl
2-propenoate (9CI)
MF (C9 H12 O6 . C6 H13 N O3 S . C4 H6 O2)x
CI PMS

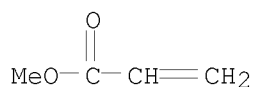
CM 1



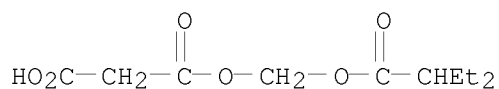
CM 2



CM 3

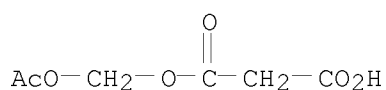


L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[(2-ethyl-1-oxobutoxy)methyl] ester
MF C10 H16 O6



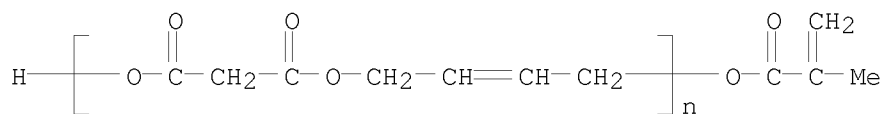
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[(acetyloxy)methyl] ester
 MF C6 H8 O6



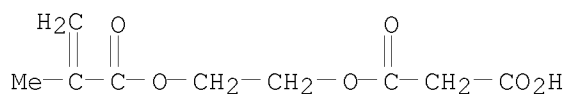
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly[oxy(1,3-dioxo-1,3-propanediyl)oxy-2-butene-1,4-diyl],
 α-hydro-ω-[(2-methyl-1-oxo-2-propenyl)oxy]- (9CI)
 MF (C7 H8 O4)_n C4 H6 O2
 CI PMS, COM

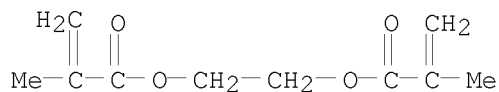


L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, mono[2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl] ester,
 polymer with 1,2-ethanediyl bis(2-methyl-2-propenoate) and methyl
 2-methyl-2-propenoate (9CI)
 MF (C10 H14 O4 . C9 H12 O6 . C5 H8 O2)_x
 CI PMS

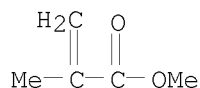
CM 1



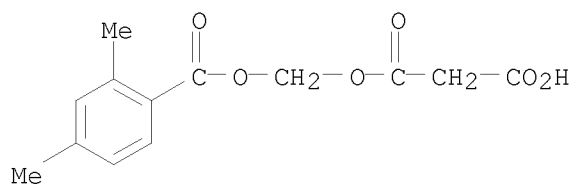
CM 2



CM 3



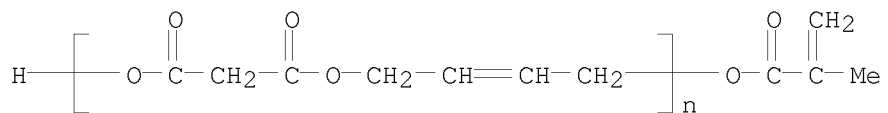
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[(2,4-dimethylbenzoyl)oxy]methyl] ester
 MF C13 H14 O6



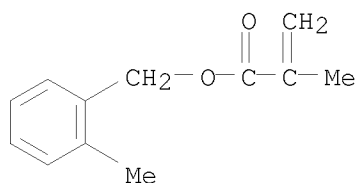
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 2-Propenoic acid, 2-methyl-, (2-methylphenyl)methyl ester, polymer with
 α -hydro- ω -[(2-methyl-1-oxo-2-propenyl)oxy]poly[oxy(1,3-dioxo-
 1,3-propanediyl)oxy-2-butene-1,4-diyl] (9CI)
 MF (C12 H14 O2 . (C7 H8 O4)_n C4 H6 O2)_x
 CI PMS

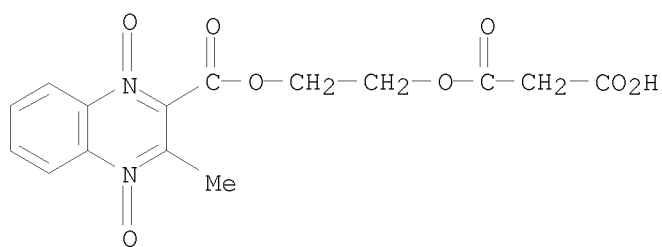
CM 1



CM 2

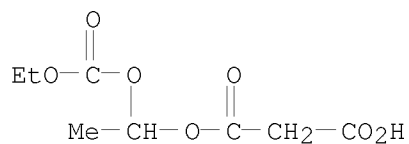


L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[[3-methyl-1,4-dioxido-2-quinoxaliny)carbonyl]oxy]ethyl] ester
 MF C15 H14 N2 O8
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

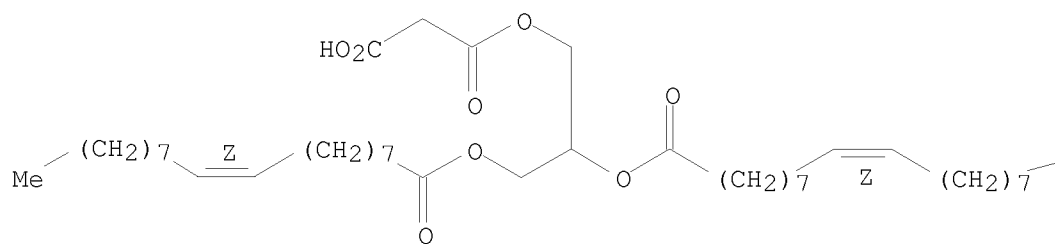
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[1-[(ethoxycarbonyl)oxy]ethyl] ester
 MF C8 H12 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2,3-bis[[9Z)-1-oxo-9-octadecen-1-yl]oxy]propyl] ester
 MF C42 H74 O8

Double bond geometry as shown.

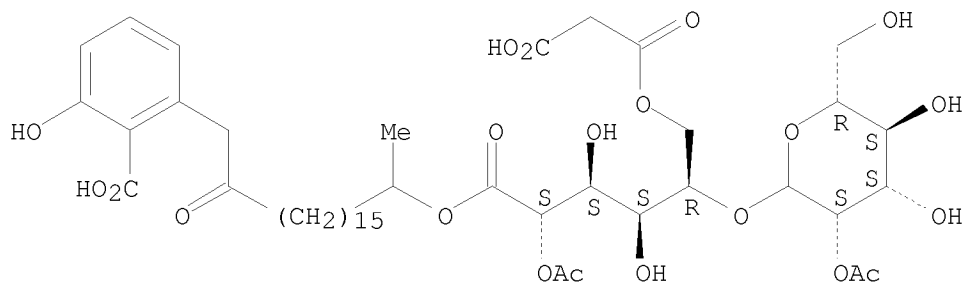


Me

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

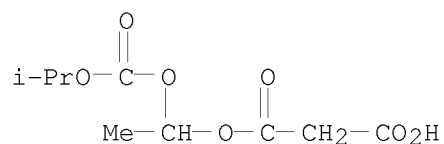
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN D-Mannonic acid, 5-O-(2-O-acetyl-D-mannopyranosyl)-,
 18-(2-carboxy-3-hydroxyphenyl)-1-methyl-17-oxooctadecyl ester, 2-acetate
 6-(hydrogen propanedioate)
 MF C45 H68 O21

Absolute stereochemistry. Rotation (-).
 Currently available stereo shown.



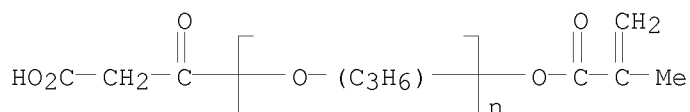
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[1-[[[(1-methylethoxy)carbonyl]oxy]ethyl] ester
 MF C9 H14 O7

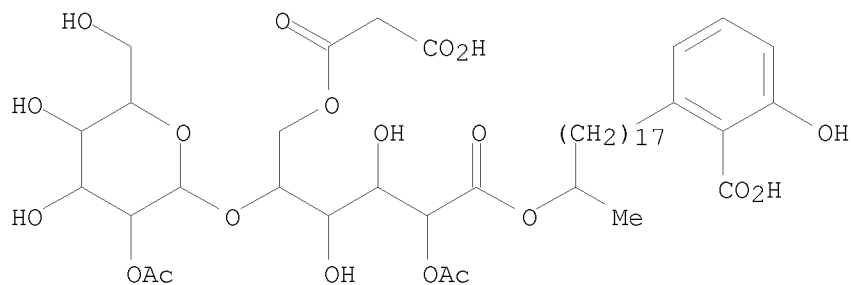


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly[oxy(methyl-1,2-ethanediyl)], α -(carboxyacetyl)- ω -[(2-methyl-1-oxo-2-propenyl)oxy]- (9CI)
 MF (C3 H6 O)_n C7 H8 O5
 CI IDS, PMS, COM

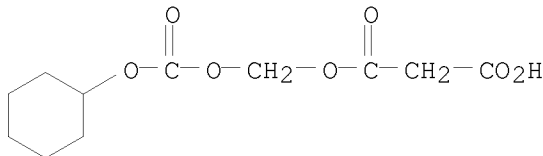


L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Hexonic acid, 5-O-(2-O-acetylhexopyranosyl)-, 18-(2-carboxy-3-hydroxyphenyl)-1-methyloctadecyl ester, 2-acetate 6-(hydrogen propanedioate) (9CI)
 MF C45 H70 O20



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

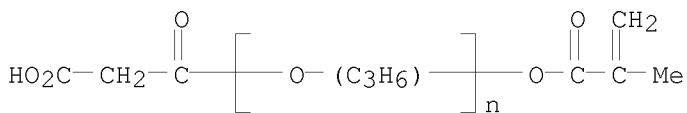
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[(cyclohexyloxy)carbonyl]oxy]methyl] ester
 MF C11 H16 O7



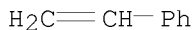
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Poly[oxy(methyl-1,2-ethanediyl)], α -(carboxyacetyl)- ω -(2-
methyl-1-oxo-2-propenyl)oxy]-, polymer with ethenylbenzene (9CI)
MF (C8 H8 . (C3 H6 O)n C7 H8 O5)x
CI PMS

CM 1

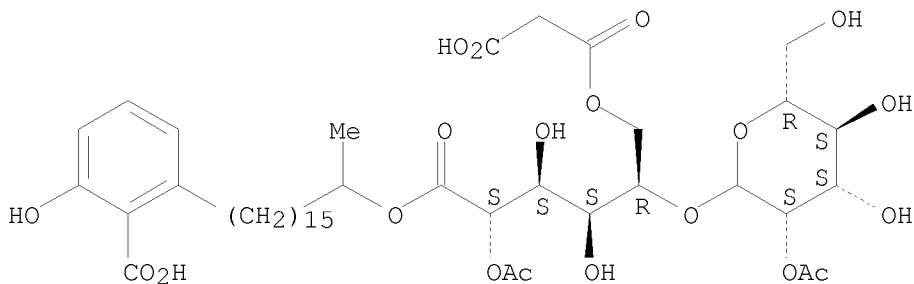


CM 2



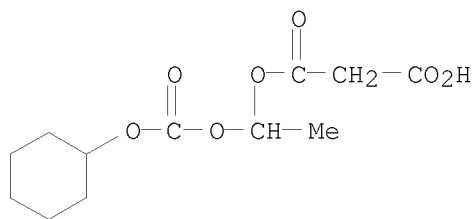
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN D-Mannonic acid, 5-O-(2-O-acetyl-D-mannopyranosyl)-,
16-(2-carboxy-3-hydroxyphenyl)-1-methylhexadecyl ester, 2-acetate
6-(hydrogen propanedioate)
MF C43 H66 O20

Absolute stereochemistry. Rotation (-).
Currently available stereo shown.



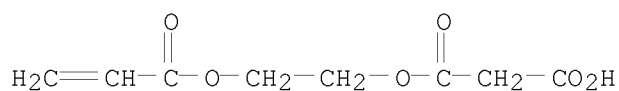
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[1-[(cyclohexyloxy)carbonyl]oxy]ethyl] ester
 MF C12 H18 O7



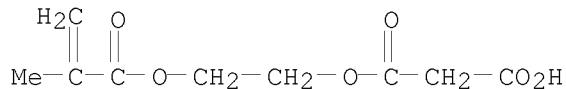
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[(1-oxo-2-propen-1-yl)oxy]ethyl] ester
 MF C8 H10 O6
 CI COM

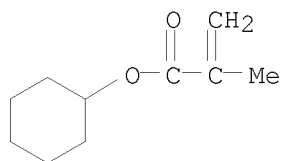


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

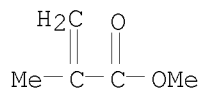
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[(2-methyl-1-oxo-2-propen-1-yl)oxy]ethyl] ester,
 polymer with cyclohexyl 2-methyl-2-propenoate and methyl
 2-methyl-2-propenoate
 MF (C10 H16 O2 . C9 H12 O6 . C5 H8 O2)x
 CI PMS
 CM 1



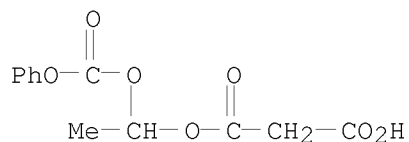
CM 2



CM 3



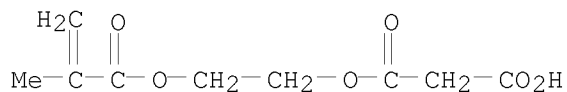
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[1-[(phoxycarbonyl)oxy]ethyl] ester
 MF C12 H12 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

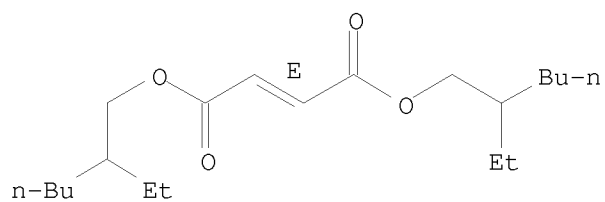
L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 2-Butenedioic acid (2E)-, bis(2-ethylhexyl) ester, polymer with
 ethenylbenzene, 2,5-furandione and 2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl
 hydrogen propanedioate (9CI)
 MF (C20 H36 O4 . C9 H12 O6 . C8 H8 . C4 H2 O3)x
 CI PMS

CM 1

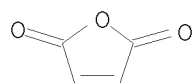


CM 2

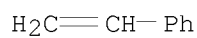
Double bond geometry as shown.



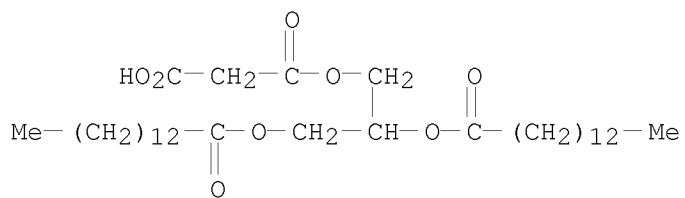
CM 3



CM 4

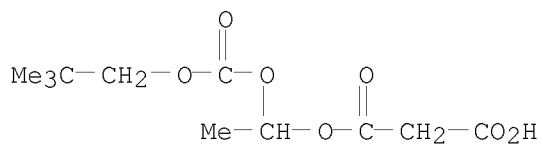


L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2,3-bis[(1-oxotetradecyl)oxy]propyl] ester
 MF C34 H62 O8



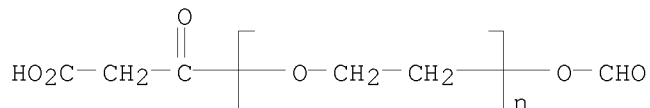
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[1-[(2,2-dimethylpropoxy)carbonyl]oxy]ethyl] ester
 MF C11 H18 O7

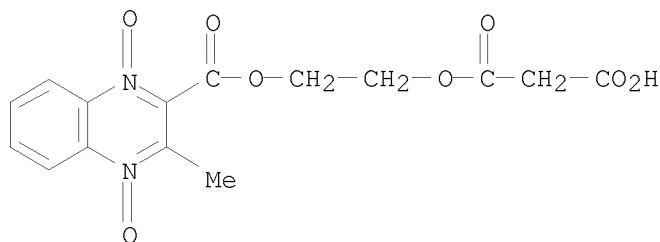


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Poly(oxy-1,2-ethanediyl), α -(carboxyacetyl)- ω -(formyloxy)-
(9CI)
MF (C2 H4 O)_n C4 H4 O5
CI PMS



L3 40 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[2-[[[3-methyl-1,4-dioxido-2-
quinoxaliny]carbonyl]oxy]ethyl] ester, sodium salt (1:1)
MF C15 H14 N2 O8 . Na



● Na

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
181.12	181.54

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 10:19:51 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available

for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 30 Dec 2008 VOL 150 ISS 1
FILE LAST UPDATED: 29 Dec 2008 (20081229/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/legal/infopolicy.html>

=> l3

L4 26 L3

=> save temp l4 alonrefs

ALONREFS IS NOT A VALID SAVED NAME

Enter the name you wish to use for the saved query, answer set, or L-number list. The name must:

1. Begin with a letter,
2. Have 1-12 characters,
3. Contain only letters (A-Z) and numbers (0-9),
4. End with /Q for a query (search profile, structure, or screen set), /A for an answer set, or /L for an L-number list.
5. Not already be in use as a saved name,
6. Not be END, SAV, SAVE, SAVED
7. Not have the form of an L-number (Lnnn).

ENTER NAME OR (END):end

=> save temp l4 malonrefs/a

ANSWER SET L4 HAS BEEN SAVED AS 'MALONREFS/A'

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	1.92	183.46

FILE 'REGISTRY' ENTERED AT 10:22:15 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7
DICTIONARY FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

Please note that search-term pricing does apply when

conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

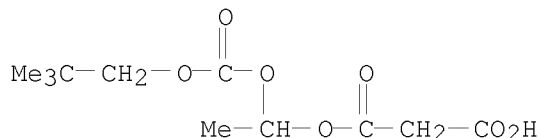
<http://www.cas.org/support/stngen/stndoc/properties.html>

```
=> e Propanedioic acid, 1-(1-(((2,2-dimethylpropoxy)carbonyl)oxy)ethyl) ester/cn
E1      1      PROPANEDIOIC ACID, 1-(1-(((1,1-DIMETHYLETHYL)AMINO)METHYL)-2
      -((9-OXO-9H-FLUOREN-4-YL)OXY)ETHYL) 3-ETHYL ESTER/CN
E2      1      PROPANEDIOIC ACID, 1-(1-(((1-METHYLETHOXY)CARBONYL)OXY)ETHYL
      ) ESTER/CN
E3      1 --> PROPANEDIOIC ACID, 1-(1-(((2,2-DIMETHYLPROPOXY)CARBONYL)OXY)
      ETHYL) ESTER/CN
E4      1      PROPANEDIOIC ACID, 1-(1-(((2-CYCLOHEXYLETHOXY)CARBONYL)OXY)E
      THYL) ESTER/CN
E5      1      PROPANEDIOIC ACID, 1-(1-(((2-METHYLPROPOXY)CARBONYL)OXY)ETHY
      L) ESTER/CN
E6      1      PROPANEDIOIC ACID, 1-(1-(((CYCLOHEXYLOXY)CARBONYL)OXY)ETHYL)
      ESTER/CN
E7      1      PROPANEDIOIC ACID, 1-(1-(((1,1-DIMETHYLETHOXY)CARBONYL)-4-PIP
      ERIDINYL) ESTER/CN
E8      1      PROPANEDIOIC ACID, 1-(1-(((3,4-DIMETHOXYPHENYL)METHYL)-2-METH
      OXY-2-OXOETHYL) ESTER/CN
E9      1      PROPANEDIOIC ACID, 1-(1-(((DIETHYLAMINO)METHYL)-2-PHENOXYETHY
      L) ESTER/CN
E10     1      PROPANEDIOIC ACID, 1-(1-(((ETHOXYCARBONYL)OXY)ETHYL) ESTER/CN
E11     1      PROPANEDIOIC ACID, 1-(1-(((PHENOXYCARBONYL)OXY)ETHYL) ESTER/C
      N
E12     1      PROPANEDIOIC ACID, 1-(1-(1,1'-BIPHENYL)-4-YLPENTYL) ESTER/CN

=> e3
L5      1 "PROPANEDIOIC ACID, 1-(1-(((2,2-DIMETHYLPROPOXY)CARBONYL)OXY)ETH
      YL) ESTER"/CN
```

=> d 15

```
L5      ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2008 ACS on STN
RN      683251-37-8  REGISTRY
ED      Entered STN:  19 May 2004
CN      Propanedioic acid, 1-[1-[[ (2,2-dimethylpropoxy)carbonyl]oxy]ethyl]
      ester (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN      Propanedioic acid, mono[1-[[ (2,2-dimethylpropoxy)carbonyl]oxy]ethyl] ester
      (9CI)
MF      C11 H18 O7
SR      CA
LC      STN Files:   CA, CAPLUS
```



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
8.99	192.45

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 10:24:34 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 30 Dec 2008 VOL 150 ISS 1
FILE LAST UPDATED: 29 Dec 2008 (20081229/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/legal/infopolicy.html>

=> 15

L6 1 L5

=> d l6 ti fbib abs

L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
TI Process for preparation of malonic acid monoesters
AN 2004:354912 CAPLUS <<LOGINID::20081230>>
DN 140:374903
TI Process for preparation of malonic acid monoesters
IN Sawabe, Takehiko; Aihara, Kazuhiro; Atsumi, Kunio; Ajito, Keiichi
PA Meiji Seika Kaisha, Ltd., Japan
SO PCT Int. Appl., 41 pp.
CODEN: PIXXD2
DT Patent
LA Japanese
FAN.CNT 2

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
-----	----	-----	-----	-----
PI WO 2004035540	A1	20040429	WO 2003-JP13319	20031017
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK,			

LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ,
 OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM,
 TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,
 KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES,
 FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
 BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

			JP 2002-304630	A	20021018
			JP 2003-50293	A	20030227
AU 2003301426	A1	20040504	AU 2003-301426		20031017
			JP 2002-304630	A	20021018
			JP 2003-50293	A	20030227
			WO 2003-JP13319	W	20031017
EP 1561748	A1	20050810	EP 2003-756680		20031017
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
			JP 2002-304630	A	20021018
			JP 2003-50293	A	20030227
			WO 2003-JP13319	W	20031017
US 20050272950	A1	20051208	US 2005-531382		20050415
			JP 2002-304630	A	20021018
			JP 2003-50293	A	20030227
			WO 2003-JP13319	W	20031017

PATENT FAMILY INFORMATION:

FAN 2004:354911

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004035539	A1	20040429	WO 2003-JP13318	20031017
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
				JP 2002-304630	A 20021018
AU 2003301425	A1	20040504	AU 2003-301425		20031017
			JP 2002-304630	A	20021018
			WO 2003-JP13318	W	20031017

OS MARPAT 140:374903

AB This invention pertains to a method for producing malonic acid monoesters with general formula of HO2CCH2CO2R [where R = a group which is easily hydrolyzed in vivo] or salts, which comprises reacting malonic acid with a halide in the presence of a base. For example, acetoxymethyl bromide was reacted with malonic acid in THF in the presence of N,N-diisopropylethylamine to give malonic acid mono-acetoxymethyl ester. This invention provides a method to make malonic acid monoesters with low cost.

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> sel 15 rn

NO ANSWERS SELECTED. THE ANSWER SET WAS CREATED IN FILE 'REGISTRY'.
 USE THE FILE COMMAND TO CHANGE TO THE CORRECT FILE.

An answer set can be processed to create terms only
 in the same file in which it was created.

=> sel l6 rn
E1 THROUGH E44 ASSIGNED

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	11.97	204.42
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.80	-0.80

FILE 'REGISTRY' ENTERED AT 10:26:51 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7
DICTIONARY FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=> s el-e44

1	103418-33-3/BI
	(103418-33-3/RN)
1	103418-34-4/BI
	(103418-34-4/RN)
1	108-48-5/BI
	(108-48-5/RN)
1	109-99-9/BI
	(109-99-9/RN)
1	1112-67-0/BI
	(1112-67-0/RN)
1	121-44-8/BI
	(121-44-8/RN)
1	141-82-2/BI
	(141-82-2/RN)
1	35180-01-9/BI
	(35180-01-9/RN)
1	40510-86-9/BI
	(40510-86-9/RN)
1	40930-71-0/BI
	(40930-71-0/RN)
1	50893-36-2/BI
	(50893-36-2/RN)
1	50972-20-8/BI

(50972-20-8/RN)
1 530-62-1/BI
(530-62-1/RN)
1 53064-79-2/BI
(53064-79-2/RN)
1 56-37-1/BI
(56-37-1/RN)
1 590-97-6/BI
(590-97-6/RN)
1 682747-70-2/BI
(682747-70-2/RN)
1 682747-74-6/BI
(682747-74-6/RN)
1 683251-13-0/BI
(683251-13-0/RN)
1 683251-19-6/BI
(683251-19-6/RN)
1 683251-21-0/BI
(683251-21-0/RN)
1 683251-24-3/BI
(683251-24-3/RN)
1 683251-28-7/BI
(683251-28-7/RN)
1 683251-31-2/BI
(683251-31-2/RN)
1 683251-33-4/BI
(683251-33-4/RN)
1 683251-34-5/BI
(683251-34-5/RN)
1 683251-37-8/BI
(683251-37-8/RN)
1 683251-39-0/BI
(683251-39-0/RN)
1 683251-42-5/BI
(683251-42-5/RN)
1 683251-45-8/BI
(683251-45-8/RN)
1 683251-48-1/BI
(683251-48-1/RN)
1 683251-50-5/BI
(683251-50-5/RN)
1 683251-53-8/BI
(683251-53-8/RN)
1 683251-62-9/BI
(683251-62-9/RN)
1 7087-68-5/BI
(7087-68-5/RN)
1 75-05-8/BI
(75-05-8/RN)
1 80715-22-6/BI
(80715-22-6/RN)
1 82504-50-5/BI
(82504-50-5/RN)
1 89838-66-4/BI
(89838-66-4/RN)
1 90776-58-2/BI
(90776-58-2/RN)
1 93457-76-2/BI
(93457-76-2/RN)
1 95775-10-3/BI
(95775-10-3/RN)

1 98298-66-9/BI
 (98298-66-9/RN)
 1 99464-83-2/BI
 (99464-83-2/RN)
 L7 44 (103418-33-3/BI OR 103418-34-4/BI OR 108-48-5/BI OR 109-99-9/BI
 OR 1112-67-0/BI OR 121-44-8/BI OR 141-82-2/BI OR 35180-01-9/BI
 OR 40510-86-9/BI OR 40930-71-0/BI OR 50893-36-2/BI OR 50972-20-8
 /BI OR 530-62-1/BI OR 53064-79-2/BI OR 56-37-1/BI OR 590-97-6/BI
 OR 682747-70-2/BI OR 682747-74-6/BI OR 683251-13-0/BI OR 683251
 -19-6/BI OR 683251-21-0/BI OR 683251-24-3/BI OR 683251-28-7/BI
 OR 683251-31-2/BI OR 683251-33-4/BI OR 683251-34-5/BI OR 683251-
 37-8/BI OR 683251-39-0/BI OR 683251-42-5/BI OR 683251-45-8/BI
 OR 683251-48-1/BI OR 683251-50-5/BI OR 683251-53-8/BI OR 683251-
 62-9/BI OR 7087-68-5/BI OR 75-05-8/BI OR 80715-22-6/BI OR 82504-
 50-5/BI OR 89838-66-4/BI OR 90776-58-2/BI OR 93457-76-2/BI OR
 95775-10-3/BI OR 98298-66-9/BI OR 99464-83-2/BI)

=> d his

(FILE 'HOME' ENTERED AT 10:14:38 ON 30 DEC 2008)

FILE 'REGISTRY' ENTERED AT 10:15:31 ON 30 DEC 2008

L1 STRUCTURE UPLOADED
 L2 2 SEARCH L1 SSS SAM
 L3 40 SEARCH L1 SSS FULL

FILE 'CAPLUS' ENTERED AT 10:19:51 ON 30 DEC 2008

L4 26 L3
 SAVE TEMP L4 MALONREFS/A

FILE 'REGISTRY' ENTERED AT 10:22:15 ON 30 DEC 2008

E PROPANEDIOIC ACID, 1-(1-((2,2-DIMETHYLPROPOXY)CARBONYL)OXY)E
 L5 1 E3

FILE 'CAPLUS' ENTERED AT 10:24:34 ON 30 DEC 2008

L6 1 L5
 SEL L6 RN

FILE 'REGISTRY' ENTERED AT 10:26:51 ON 30 DEC 2008

L7 44 S E1-E44

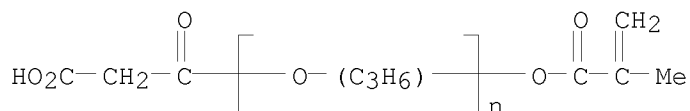
=> l3 not l7

L8 25 L3 NOT L7

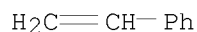
=> d scan

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly[oxy(methyl-1,2-ethanediyl)], α -(carboxyacetyl)- ω -[(2-
 methyl-1-oxo-2-propenyl)oxy]-, polymer with ethenylbenzene (9CI)
 MF (C8 H8 . (C3 H6 O)n C7 H8 O5)x
 CI PMS

CM 1



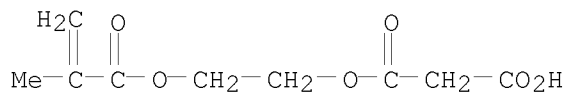
CM 2



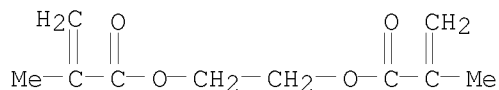
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):25

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, mono[2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl] ester,
polymer with 1,2-ethanediyl bis(2-methyl-2-propenoate) and methyl
2-methyl-2-propenoate (9CI)
MF (C10 H14 O4 . C9 H12 O6 . C5 H8 O2)x
CI PMS

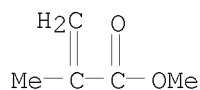
CM 1



CM 2

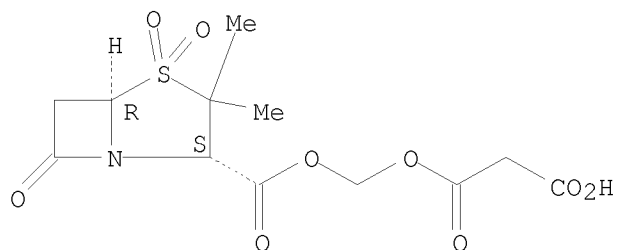


CM 3



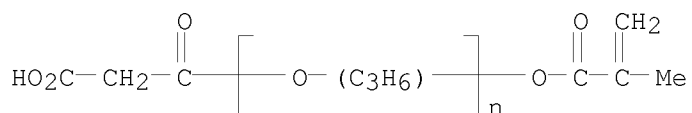
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-
azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, sodium salt,
(2S-cis)- (9CI)
MF C12 H15 N O9 S . Na

Absolute stereochemistry.



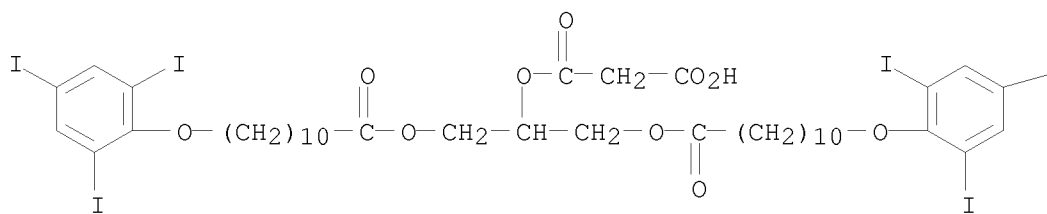
● Na

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly[oxy(methyl-1,2-ethanediyl)], α -(carboxyacetyl)- ω -(2-methyl-1-oxo-2-propenyl)oxy]- (9CI)
 MF (C3 H6 O)_n C7 H8 O5
 CI IDS, PMS, COM



L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[[1-oxo-11-(2,4,6-triiodophenoxy)undecyl]oxy]-1-[[[1-oxo-11-(2,4,6-triiodophenoxy)undecyl]oxy]methyl]ethyl] ester
 MF C40 H52 I6 O10

PAGE 1-A

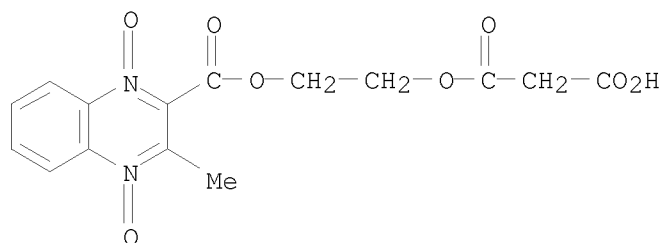


PAGE 1-B

— I

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[[[(3-methyl-1,4-dioxido-2-
 quinoxaliny)carbonyl]oxy]ethyl] ester, sodium salt (1:1)
 MF C15 H14 N2 O8 . Na

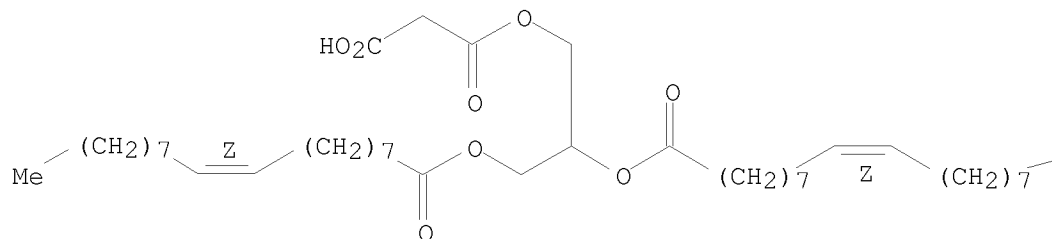


● Na

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2,3-bis[[[(9Z)-1-oxo-9-octadecen-1-yl]oxy]propyl]
 ester
 MF C42 H74 O8

Double bond geometry as shown.

PAGE 1-A

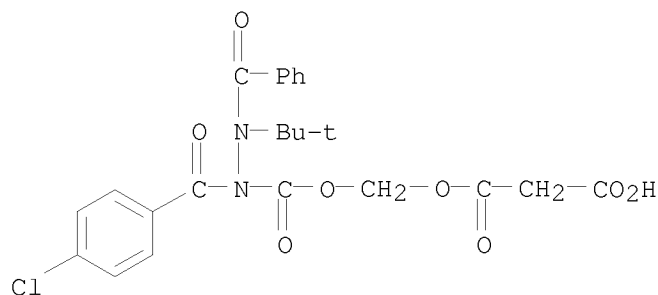


PAGE 1-B

Me

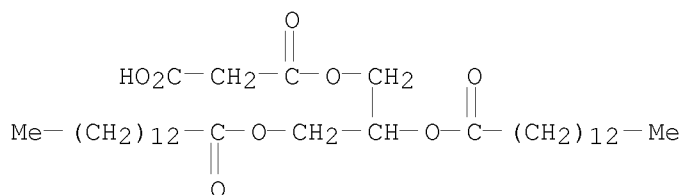
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[2-benzoyl-1-(4-chlorobenzoyl)-2-(1,1-dimethylethyl)hydrazinyl]carbonyl]oxy]methyl] ester, sodium salt (1:1)
 MF C23 H23 Cl N2 O8 . Na



● Na

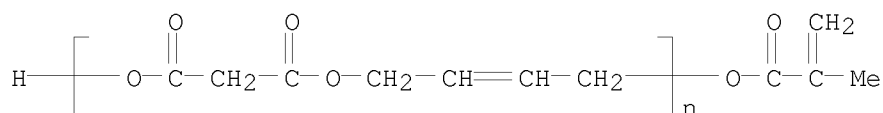
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2,3-bis[(1-oxotetradecyl)oxy]propyl] ester
 MF C34 H62 O8



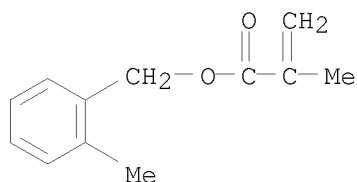
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 2-Propenoic acid, 2-methyl-, (2-methylphenyl)methyl ester, polymer with α -hydro- ω -[(2-methyl-1-oxo-2-propenyl)oxy]poly[oxy(1,3-dioxo-1,3-propanediyl)oxy-2-butene-1,4-diyl] (9CI)
 MF (C12 H14 O2 . (C7 H8 O4)n C4 H6 O2)x
 CI PMS

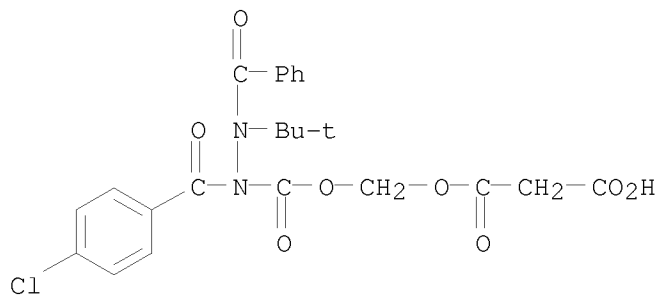
CM 1



CM 2



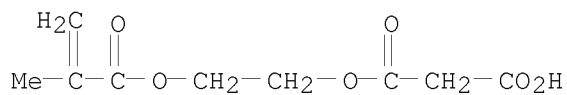
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[[[2-benzoyl-1-(4-chlorobenzoyl)-2-(1,1-dimethylethyl)hydrazinyl]carbonyl]oxy]methyl] ester
 MF C23 H23 Cl N2 O8
 CI COM



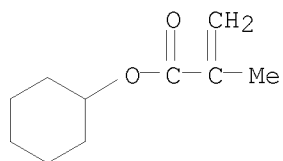
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[(2-methyl-1-oxo-2-propen-1-yl)oxy]ethyl] ester,
 polymer with cyclohexyl 2-methyl-2-propenoate and methyl
 2-methyl-2-propenoate
 MF (C10 H16 O2 . C9 H12 O6 . C5 H8 O2)x
 CI PMS

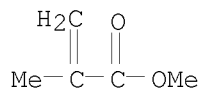
CM 1



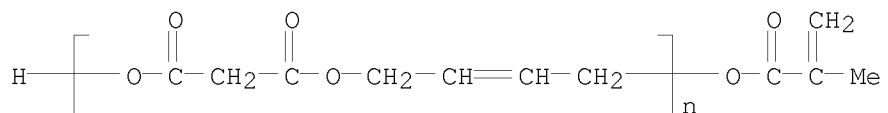
CM 2



CM 3

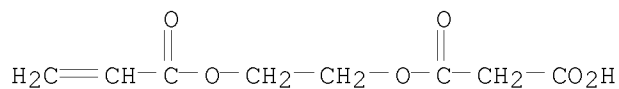


L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly[oxy(1,3-dioxo-1,3-propanediyl)oxy-2-butene-1,4-diyl],
 α -hydro- ω -[(2-methyl-1-oxo-2-propenyl)oxy]- (9CI)
 MF (C7 H8 O4)_n C4 H6 O2
 CI PMS, COM



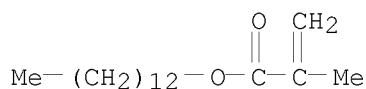
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 2-Propenoic acid, 2-methyl-, tridecyl ester, polymer with dodecyl
 2-propenoate and 2-hydroxyethyl 2-propenoate,
 2-[(1-oxo-2-propenyl)oxy]ethyl propanedioate (9CI)
 MF (C17 H32 O2 . C15 H28 O2 . C5 H8 O3)_x . x C8 H10 O6

CM 1

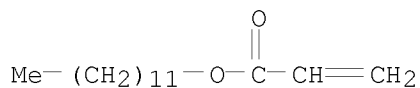


CM 2

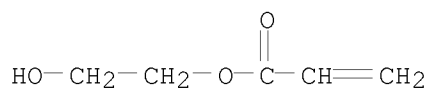
CM 3



CM 4

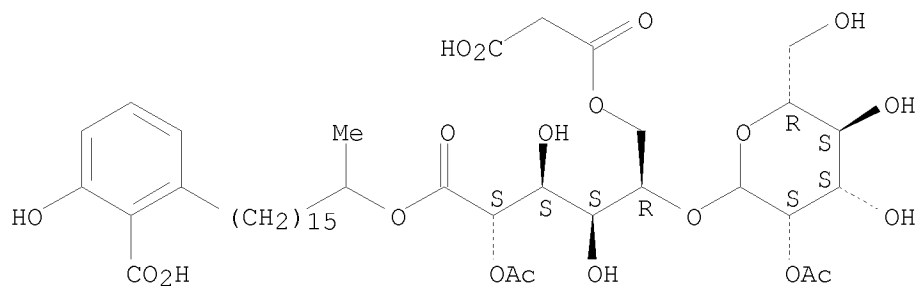


CM 5



L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN D-Mannonic acid, 5-O-(2-O-acetyl-D-mannopyranosyl)-,
16-(2-carboxy-3-hydroxyphenyl)-1-methylhexadecyl ester, 2-acetate
6-(hydrogen propanedioate)
MF C43 H66 O20

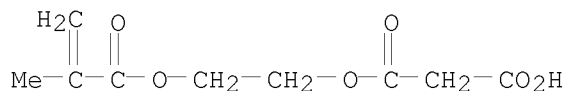
Absolute stereochemistry. Rotation (-).
Currently available stereo shown.



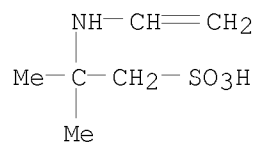
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, mono[2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl] ester,
polymer with 2-(ethenylamino)-2-methyl-1-propanesulfonic acid and methyl
2-propenoate (9CI)
MF (C9 H12 O6 . C6 H13 N O3 S . C4 H6 O2)x
CI PMS

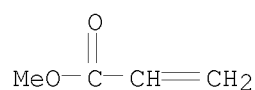
CM 1



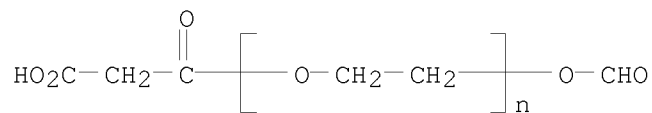
CM 2



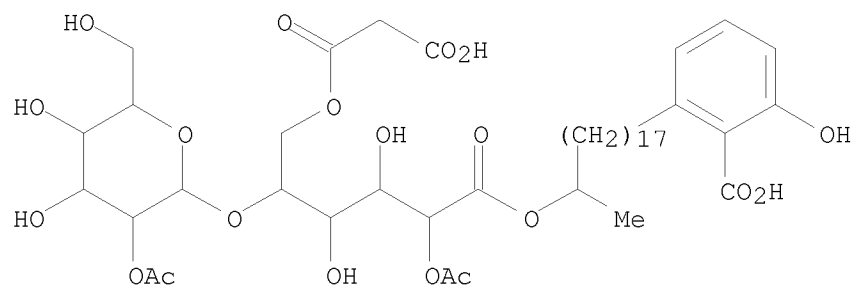
CM 3



L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Poly(oxy-1,2-ethanediyl), α -(carboxyacetyl)- ω -(formyloxy)-
 (9CI)
 MF (C2 H4 O)_n C4 H4 O5
 CI PMS

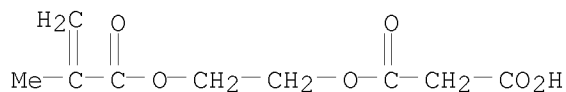


L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Hexonic acid, 5-O-(2-O-acetylhexopyranosyl)-,
 18-(2-carboxy-3-hydroxyphenyl)-1-methyloctadecyl ester, 2-acetate
 6-(hydrogen propanedioate) (9CI)
 MF C45 H70 O20



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

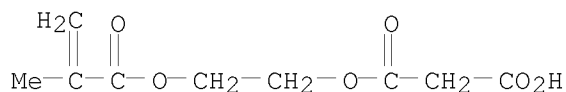
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[(2-methyl-1-oxo-2-propen-1-yl)oxy]ethyl] ester
 MF C9 H12 O6
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

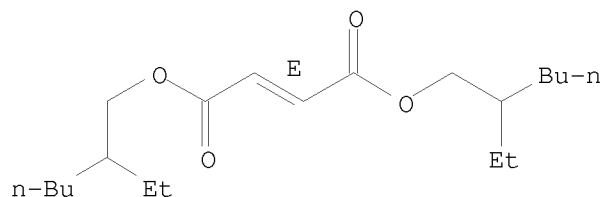
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 2-Butenedioic acid (2E)-, bis(2-ethylhexyl) ester, polymer with
 ethenylbenzene, 2,5-furandione and 2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl
 hydrogen propanedioate (9CI)
 MF (C20 H36 O4 . C9 H12 O6 . C8 H8 . C4 H2 O3)x
 CI PMS

CM 1

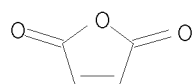


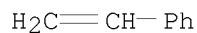
CM 2

Double bond geometry as shown.



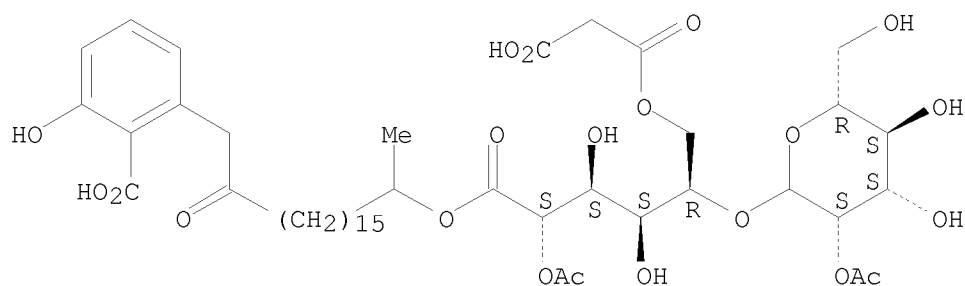
CM 3





L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN D-Mannonic acid, 5-O-(2-O-acetyl-D-mannopyranosyl)-,
 18-(2-carboxy-3-hydroxyphenyl)-1-methyl-17-oxooctadecyl ester, 2-acetate
 6-(hydrogen propanedioate)
 MF C45 H68 O21

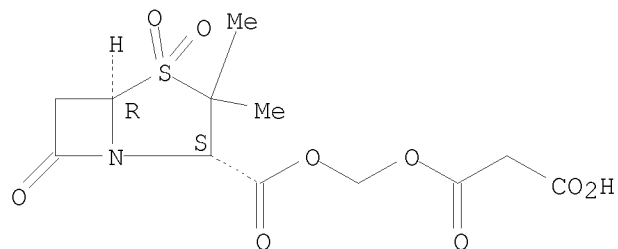
Absolute stereochemistry. Rotation (-).
 Currently available stereo shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, mono[[[(3,3-dimethyl-4,4-dioxido-7-oxo-4-thia-1-azabicyclo[3.2.0]hept-2-yl)carbonyl]oxy]methyl] ester, (2S-cis)- (9CI)
 MF C12 H15 N O9 S
 CI COM

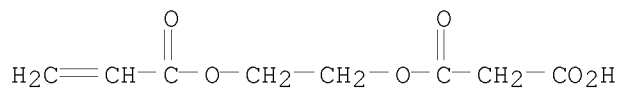
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

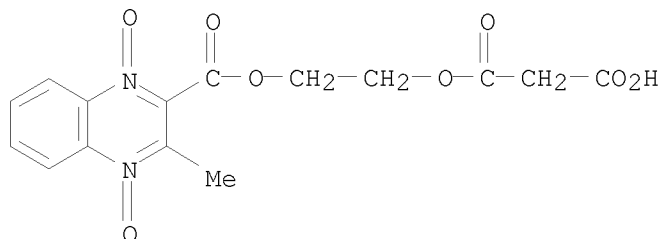
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Propanedioic acid, 1-[2-[(1-oxo-2-propen-1-yl)oxy]ethyl] ester

MF C8 H10 O6
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

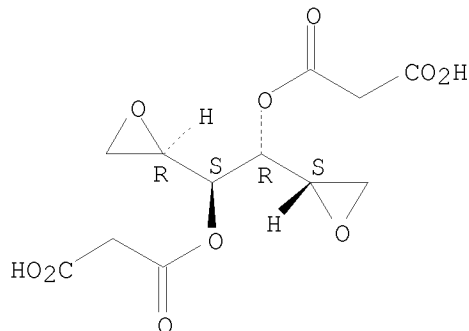
L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Propanedioic acid, 1-[2-[[3-methyl-1,4-dioxido-2-
quinoxaliny]carbonyl]oxy]ethyl] ester
MF C15 H14 N2 O8
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 25 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Galactitol, 1,2:5,6-dianhydro-, bis(hydrogen propanedioate) (9CI)
MF C12 H14 O10

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.22	207.64
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.80

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 10:30:58 ON 30 DEC 2008

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'REGISTRY' AT 11:24:17 ON 30 DEC 2008
FILE 'REGISTRY' ENTERED AT 11:24:17 ON 30 DEC 2008
COPYRIGHT (C) 2008 American Chemical Society (ACS)

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.22	207.64
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.80

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.68	208.10
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.80

FILE 'REGISTRY' ENTERED AT 11:25:07 ON 30 DEC 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7
DICTIONARY FILE UPDATES: 29 DEC 2008 HIGHEST RN 1091682-77-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary files\10531382\10531382 phthalidyl.str

L9 STRUCTURE UPLOADED

=> d 15

L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2008 ACS on STN

RN 683251-37-8 REGISTRY

ED Entered STN: 19 May 2004

CN Propanedioic acid, 1-[1-[[(2,2-dimethylpropoxy)carbonyl]oxy]ethyl] ester (CA INDEX NAME)

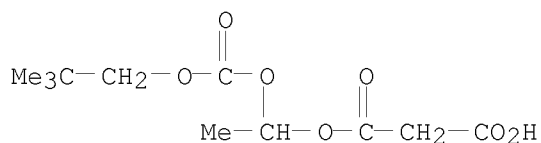
OTHER CA INDEX NAMES:

CN Propanedioic acid, mono[1-[[(2,2-dimethylpropoxy)carbonyl]oxy]ethyl] ester (9CI)

MF C11 H18 O7

SR CA

LC STN Files: CA, CAPLUS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

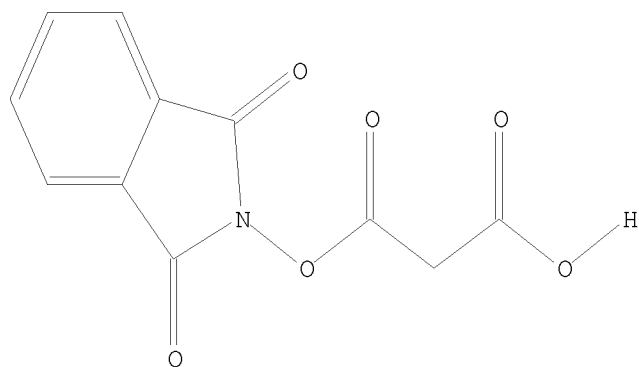
1 REFERENCES IN FILE CA (1907 TO DATE)

1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d 19

L9 HAS NO ANSWERS

L9 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l9 sss sam

SAMPLE SEARCH INITIATED 11:25:58 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 83 TO ITERATE

100.0% PROCESSED 83 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 1114 TO 2206

PROJECTED ANSWERS: 0 TO 0

L10 0 SEA SSS SAM L9

=> search l9 sss full

FULL SEARCH INITIATED 11:26:06 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 1151 TO ITERATE

100.0% PROCESSED 1151 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

L11 0 SEA SSS FUL L9

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	185.88	393.98
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.80

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 11:32:47 ON 30 DEC 2008